EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	2	("5739163").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR ,	OFF	2008/01/23 05:48
L2	. 2	("5523302").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2008/01/23 05:48
L3	6	"9736862"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR .	ON	2008/01/23 05:48
L4	2	("5773646").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2008/01/23 05:48
L5	18	("5523302").URPN.	USPAT	OR	ON	2008/01/23 05:48
L6	354	(562/431).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2008/01/23 05:48
L7	386	(562/471).CCLS.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2008/01/23 05:48
L8	317	(562/472).CCLS. ¹	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	OFF	2008/01/23 05:48
L9	910	L6 or L7 or L8	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48
L10	28904	williamson	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48
L11	11	L9 and L10	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48
L12	0	"4656305.pn"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48

EAST Search History

L13	0	"4656305.pn."	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48
L14	2.	"4656305".pn.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:48
L15	2	("4150238").PN.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR .	OFF	2008/01/23 05:52
L16	2	"4656305".pn.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 05:53
L18	8	"1,3-diphenylprop-2-en-1-one"	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 06:28
L19	4	"1,3-diphenylprop-2-en-1-one".clm.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2008/01/23 06:32

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        OCT 19
                 BEILSTEIN updated with new compounds
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        NOV 15
NEWS 17
                 WPIX enhanced with XML display format
NEWS 18 NOV 19
NEWS 19 NOV 30
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NEWS 20 DEC 04
                 LINPADOCDB now available on STN
NEWS 21 DEC 14
                 BEILSTEIN pricing structure to change
NEWS 22 DEC 17
                 USPATOLD added to additional database clusters
                 IMSDRUGCONF removed from database clusters and STN
NEWS 23 DEC 17
NEWS 24 DEC 17
                 DGENE now includes more than 10 million sequences
NEWS 25 DEC 17
                 TOXCENTER enhanced with 2008 MeSH vocabulary in
                 MEDLINE segment
NEWS 26 DEC 17
                 MEDLINE and LMEDLINE updated with 2008 MeSH vocabulary
NEWS 27 DEC 17
                 CA/CAplus enhanced with new custom IPC display formats
                 STN Viewer enhanced with full-text patent content
NEWS 28
       DEC 17
                 from USPATOLD
NEWS 29
         JAN 02
                 STN pricing information for 2008 now available
NEWS 30
        JAN 16
                 CAS patent coverage enhanced to include exemplified
                 prophetic substances
NEWS EXPRESS
              19 SEPTEMBER 2007: CURRENT WINDOWS VERSION IS V8.2,
              CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
              AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.
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=> file reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 1.47 1.47

FULL ESTIMATED COST

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L1 STRUCTURE UPLOADED

=> d l1

≃>

L1 HAS NO ANSWERS

L1 STR

Structure attributes must be viewed using STN Express query preparation.

=> .

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L2 STRUCTURE UPLOADED

=> d 12

L2 HAS NO ANSWERS

L2 STR

Structure attributes must be viewed using STN Express query preparation.

=> search 12 sss sam

SAMPLE SEARCH INITIATED 08:11:08 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 76 TO ITERATE

100.0% PROCESSED 76 ITERATIONS 2 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 997 TO 2043

PROJECTED ANSWERS: 2 TO 124

L3 2 SEA SSS SAM L2

=> d scan

L3 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN 1,4-Benzenedicarboxylic acid, dimethyl ester, polymer with α -[[4-(carboxymethoxy)phenyl]methylene]- β -oxobenzenepropanoic acid, 1,4-cyclohexanedimethanol and 1,2-ethanediol (9CI)

MF (C18 H14 O6 . C10 H10 O4 . C8 H16 O2 . C2 H6 O2)x

CI PMS

CM 1

$$\begin{array}{c|c} & \text{HO}_2\text{C} & \text{O} \\ & \parallel \\ \text{CH} & \text{C-C-Ph} \\ \\ \text{HO}_2\text{C-CH}_2 - \text{O} \end{array}$$

CM 2

CM 3

 ${\tt HO-CH_2-CH_2-OH}$

CM 4

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Acetic acid, [2-(3-oxo-3-phenyl-1-propenyl)phenoxy]-, (E)- (9CI)
MF C17 H14 O4

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

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=> 13

L4 3 L3

=> d 14 1-3 ti fbib abs

L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN

TI Aryloxymethyl radical cyclizations mimicking biological carbon-carbon bond formation to methoxy groups

AN 1993:147350 CAPLUS <<LOGINID::20080123>>

DN 118:147350

TI Aryloxymethyl radical cyclizations mimicking biological carbon-carbon bond formation to methoxy groups

AU Ahmad-Junan, S. Asiah; Walkington, Andrew J.; Whiting, Donald A.

CS Dep. Chem., Univ. Nottingham, Nottingham, NG7 2RD, UK

SO Journal of the Chemical Society, Perkin Transactions 1: Organic and Bio-Organic Chemistry (1972-1999) (1992), (18), 2313-20 CODEN: JCPRB4; ISSN: 0300-922X

DT Journal

LA English

OS CASREACT 118:147350

GI

As small group of diverse natural products whose biosynthesis is unusual in involving formation of O-heterocyclic rings by C-C bond formation to aromatic methoxy groups, in net oxidative and non-oxidative processes is examined It is shown that aryloxymethyl radicals, generated by decarboxylation of thiohydroxamate esters of aryloxyacetic acids, undergo addition to both electron-rich and electron-poor double bonds, cyclizing in 5-exo, 6-endo and 6-exo fashions, and also substitute regioselectively into pyridinium nuclei, thus mimicking the biochem. processes, as well as forming a useful new synthetic approach to O heterocycles. In an example the isoderritol isoflavone I is cyclized to dehydrorotenone via I (R = CH2CO2H). Biosynthetic mechanisms are briefly discussed.

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN

TI Biomimetic radical cyclizations: synthesis and biosynthesis of benzodihydropyrans and -furans

AN 1990:198051 CAPLUS <<LOGINID::20080123>>

DN 112:198051

TI Biomimetic radical cyclizations: synthesis and biosynthesis of benzodihydropyrans and -furans

AU Walkington, Andrew J.; Whiting, Donald A.

CS Chem. Dep., Univ. Nottingham, Nottingham, NG7 2RD, UK

SO Tetrahedron Letters (1989), 30(35), 4731-4

CODEN: TELEAY; ISSN: 0040-4039

DT Journal

LA English

OS CASREACT 112:198051

GI

AB Aryloxymethylene radicals generated by decarboxylaton of the thiohydroxamate esters derived from acids 2,6-RR1C6H3OCH2CO2H R = alkyl, R1 = OMe; R = COCH:CHPh, R1 = H) and 1,2-(Me2C:CHCH2)C10H6OCH2CO2H undergo 6-endo cyclization yielding I-III resp., mimicking the unusual biosynthetic reactions involved in formation of scabequinone, benzylchromanone, stachyoidin in nature.

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN

TI Stabilized copolyester material

AN 1982:564084 CAPLUS <<LOGINID::20080123>>

DN 97:164084

OREF 97:27385a,27388a

TI Stabilized copolyester material

IN Zannucci, Joseph S.; Sublett, Bobby J.; Irick, Gether, Jr.

PA Eastman Kodak Co., USA

SO U.S., 5 pp.

CODEN: USXXAM

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ΡI	US 4340718	A	19820720	US 1980-155811	19800602
				US 1980-155811	19800602

GI

Polyesters having good weatherability contain 0.3-5.0% stabilizing moieties I (R = 1-4 groups independently selected from H, Cl, F, alkyl, cycloalkyl, alkoxy, aryl, or aryloxy; R1 = alkyl, cycloalkyl, or aryl; R2 = CN, alkylcarbonyl, cycloalkylcarbonyl, or arylcarbonyls; Z = CO, O2CR3O, CO2R3O, O2CR3, O2CO, or O2COR3O, or R3 = C1-10 alkylene, arylene, or cycloalkylene). Thus, a polyester [83266-00-6] prepared from di-Me terephthalate, ethylene glycol, 1,4-cyclohexanedimethanol, and p-MeO2CCH2OC6H4CH:C(CN)CO2Me (1.0% of total polymer moiety) had impact strength after 500 h exposure to a weather ometer ≥7 + that of the unstabilized polyester.

=> file reg COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	12.09	16.78
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-2.40	2.40

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=>

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L5 STRUCTURE UPLOADED

=> d 15

L5 HAS NO ANSWERS

L5 STR

Structure attributes must be viewed using STN Express query preparation.

=> search 15 sss sam

SAMPLE SEARCH INITIATED 08:16:23 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 76 TO ITERATE

100.0% PROCESSED 76 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 997 TO 2043

PROJECTED ANSWERS: 1 TO 80

L6 1 SEA SSS SAM L5

=> d scan

L6 1 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Acetic acid, [2-(3-oxo-3-phenyl-1-propenyl)phenoxy]-, (E)- (9CI)

MF C17 H14 O4

Double bond geometry as shown.

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL. ENTRY SESSION 0.46 17.24 FULL ESTIMATED COST DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL SESSION ENTRY -2.40 CA SUBSCRIBER PRICE 0.00

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=> 16

L7 2 L6

=> d 17 1-2 ti

- L7 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN
- TI Aryloxymethyl radical cyclizations mimicking biological carbon-carbon bond formation to methoxy groups
- L7 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN
- TI Biomimetic radical cyclizations: synthesis and biosynthesis of benzodihydropyrans and -furans

=> file reg

COST IN U.S. DOLLARS SINCE FILE TOTAL SESSION ENTRY FULL ESTIMATED COST 8.88 26.12 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -2.40

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=> d his

(FILE 'HOME' ENTERED AT 08:03:33 ON 23 JAN 2008)

FILE 'REGISTRY' ENTERED AT 08:07:54 ON 23 JAN 2008

L1 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED
L3 2 SEARCH L2 SSS SAM

FILE 'CAPLUS' ENTERED AT 08:12:03 ON 23 JAN 2008

L4 3 L3

FILE 'REGISTRY' ENTERED AT 08:15:58 ON 23 JAN 2008

L5 STRUCTURE UPLOADED

L6 1 SEARCH L5 SSS SAM

FILE 'CAPLUS' ENTERED AT 08:16:38 ON 23 JAN 2008 L7 2 L6

FILE 'REGISTRY' ENTERED AT 08:26:55 ON 23 JAN 2008

=> search 15 sss full

FULL SEARCH INITIATED 08:27:22 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1369 TO ITERATE

100.0% PROCESSED 1369 ITERATIONS

6 ANSWERS

SEARCH TIME: 00.00.01

L8 6 SEA SSS FUL L5

=> d scan

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Acetic acid, [2,3-dichloro-4-(3-oxo-3-phenyl-1-propenyl)phenoxy]- (9CI)

MF C17 H12 C12 O4

$$CH = CH - C - Ph$$
 $C1$

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):6

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Acetic acid, [2-(3-oxo-3-phenyl-1-propenyl)phenoxy]-, (E)- (9CI)
MF C17 H14 O4

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Acetic acid, [4-benzo[b]thien-2-yl-5-methoxy-2-[(1E)-3-oxo-3-phenyl-1-propenyl]phenoxy]- (9CI)
MF C26 H20 O5 S

Double bond geometry as shown.

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN IN Acetic acid, [m-(2-benzoylvinyl)phenoxy]- (5CI) MF C17 H14 O4

$$HO_2C-CH_2-O$$
 CH
 CH
 CH
 CH

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Acetic acid, 2-[2-methoxy-4-[(1Z)-3-oxo-3-phenyl-1-propen-1-yl]phenoxy]MF C18 H16 O5

Double bond geometry as shown.

$$MeO$$
 Z
 O
 Ph

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 6 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Acetic acid, [4-(3-oxo-3-phenyl-1-propenyl)phenoxy]- (9CI)
MF C17 H14 O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> e Acetic acid, (4-(3-oxo-3-phenyl-1-propenyl)phenoxy)-/cn

```
ACETIC ACID, (4-(3-OXO-3-(4-SULFOPHENYL)-1-PROPENYL) PHENOXY)
E1
                   -, (E) - /CN
                   ACETIC ACID, (4-(3-OXO-3-(4-SULFOPHENYL)-1-PROPENYL) PHENOXY)
E2
                   -, DISODIUM SALT, (E)-/CN
             1 --> ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL)PHENOXY)-/CN
E3
                   ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL)PHENOXY)-, ETHYL
E4
                   ESTER/CN
                   ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL) PHENOXY)-, ETHYL
E5
                   ESTER, (E)-/CN
                   ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL)PHENOXY)-, METHYL
E6
                    ESTER/CN
                   ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL) PHENOXY)-, METHYL
E7
                    ESTER, (E)-/CN
                   ACETIC ACID, (4-(3-OXO-6-(4-PHENYL-1-BUTENYL)BICYCLO(3.1.0)H
E8
                   EX-2-YL)BUTOXY)-, ETHYL ESTER, (1A,2A,5A,6
                   B(Z))-/CN
                   ACETIC ACID, (4-(3-OXO-6-(4-PHENYL-1-BUTENYL)BICYCLO(3.1.0)H
             1
E9
                   EX-2-YL)BUTOXY)-, ETHYL ESTER, (1A, 2B, 5A, 6.
                   BETA. (Z)) - /CN
                   ACETIC ACID, (4-(3-OXO-6-(4-PHENYL-1-BUTENYL)BICYCLO(3.1.0)H
             1
E10
                   EX-2-YL) BUTOXY) -, ETHYL ESTER, (1A, 2B, 5A, 6.
                   BETA. (Z)) - (\pm) - /CN
                   ACETIC ACID, (4-(3-OXO-6-(4-PHENYL-1-BUTENYL)BICYCLO(3.1.0)H
E11
             1
                   EX-2-YL)BUTOXY)-, ETHYL ESTER, (Z)-/CN
                   ACETIC ACID, (4-(3-OXO-6-(4-PHENYL-1-BUTENYL)BICYCLO(3.1.0)H
E12
             1
                   EX-2A-YL) BUTOXY) -, ETHYL ESTER, (Z)-ENDO-/CN
=> e3
             1 "ACETIC ACID, (4-(3-OXO-3-PHENYL-1-PROPENYL) PHENOXY) - "/CN
L9
=> d 19
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2008 ACS on STN
L9
RN
     31824-93-8 REGISTRY
ED
     Entered STN: 16 Nov 1984
     Acetic acid, [4-(3-oxo-3-phenyl-1-propenyl)phenoxy]- (9CI)
CN
     INDEX NAME)
OTHER CA INDEX NAMES:
     Acetic acid, [p-(2-benzoylvinyl)phenoxy] - (8CI)
CN
MF
     C17 H14 O4
                  BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMCATS, IFICDB, IFIPAT,
LC
       IFIUDB, TOXCENTER, USPATOLD
```

$$CH = CH - C - Ph$$

$$HO_2C - CH_2 - O$$

- 4 REFERENCES IN FILE CA (1907 TO DATE)
- 4 REFERENCES IN FILE CAPLUS (1907 TO DATE)

(*File contains numerically searchable property data)

=> file caplus
COST IN U.S. DOLLARS

FULL ESTIMATED COST ENTRY SESSION 186.89 213.01

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
SINCE FILE TOTAL ENTRY SESSION

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=> 19

L10 4 L9

=> d 110 1-4 ti fbib abs

L10 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN

- TI 4-Substituted phenoxyacetic acid derivatives and their antimicrobial activities
- AN 1992:193834 CAPLUS <<LOGINID::20080123>>
- DN 116:193834
- TI 4-Substituted phenoxyacetic acid derivatives and their antimicrobial activities
- AU Sarac, Selma; Safak, Cihat; Erdogan, Hakki; Abbasoglu, Ufuk; Gunay, Yekta
- CS Fac. Pharm., Hacettepe Univ., Ankara, Turk.
- SO Hacettepe Universitesi Eczacilik Fakultesi Dergisi (1991), 11(1), 1-11 CODEN: HUEDEE: ISSN: 1300-0608
- DT Journal
- LA English
- OS CASREACT 116:193834

GI

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2-thienyl) were prepared by condensing phenoxyacetate II (R1 = CHO) with
    RCOMe in the presence of NaOH in EtOH. Title compds. III (R = same) were
    prepared by condensing II (R1 = Ac) with aldehydes RCHO in the presence of
    NaOH in EtOH. I and III exhibited antifungal and antibacterial
    activities. ·
    ANSWER 2 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
TI
    Substituted phenoxyacetic acids
    1974:145808 CAPLUS <<LOGINID::20080123>>
AN
     80:145808
DN
OREF 80:23525a,23528a
    Substituted phenoxyacetic acids
TI
    Kuhn, Stephen J.; Ilavsky, Janet E.
IN
PΑ
    Dow Chemical Co.
SO
    Can., 9 pp.
    CODEN: CAXXA4
DT
    Patent
LA
    English
FAN.CNT 1
                       KIND
                                          APPLICATION NO.
                                                                 DATE
    PATENT NO.
                               DATE
                                           ______
                               _____
     _____
                                         CA 1970-93423
                               19720704
                                                                 19700917
    CA 904291
PI
    p-RCOCH:CHC6H4OCH2CO2H (R = p-BrC6H4, p-EtOC6H4, 1-C10H7, Ph, p-ClC6H4,
AB
    p-PrC6H4, 3,4-Cl2C6H3, 2,4-Et2C6H3, p-EtC6H4, 2,4-ClEtC6H3, 2-thienyl,
     2,4-Br(Me2CHO)C6H3, 5-indanyl), useful as pesticides, were prepared by the
     condensation of RCOMe with p-OCHC6H4OCH2CO2H.
L10 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
    Pesticidal substituted phenoxyacetic acids
TI
     1971:125177 CAPLUS <<LOGINID::20080123>>
AN
DN
    74:125177
OREF 74:20219a,20222a
    Pesticidal substituted phenoxyacetic acids
     Kuhn, Stephen J.; Ilavsky, Janet E.
    Dow Chemical Co.
PA
    U.S., 3 pp.
SO
    CODEN: USXXAM
DT
    Patent
LΑ
    English
FAN.CNT 1
                                          APPLICATION NO.
                                                                 DATE
     PATENT NO.
                        KIND
                               DATE
                                          ______
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                               _____
                                          US 1968-731292
                                                                  19680522
                               19710126
PΙ
    US 3558612
                                                             A 19680522
                                           US 1968-731292
     For diagram(s), see printed CA Issue.
GI
     The title compds. (I) useful as pesticides, are prepared by condensation
AB
     from 4-formylphenoxyacetic acid and a p-substituted acetophenone in aqueous
     alc. or by substitution of a p-bromo compound and a suitable sodio derivative
by
     metathesis. The following I were prepared (X and Y given): 4-Br, H; 4-EtO,
     H; H, H; 4-Cl, H; 3-Cl, 4-Cl; 4-Et, H; 4-EtO, H. The 4-(2-
     naphthylcarbonylvinyl), the 4-(2-thienylcarbonylvinyl) and the
     4-(5-indanylcarbonylvinyl) derivs. were also prepared
    ANSWER 4 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
L10
     Chemotherapeutic thiosemicarbazones
TT
     1956:4949 CAPLUS <<LOGINID::20080123>>
AN
DN
     50:4949
OREF 50:1086a-d
     Chemotherapeutic thiosemicarbazones
TI
     Farbenfabriken Bayer A.-G.
PA
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Title compds. I (R = Ph, 4-MeC6H4, 4-MeOC6H4, 4-ClC6H4, 4-BrC6H4,

AB

DT

Patent

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Unavailable
LA
FAN.CNT 1
                                 DATE
                                             APPLICATION NO.
     PATENT NO.
                          KIND
                                 19540428
                                            GB
ΡI
     GB 708013
     Substituted styryl aryl ketone thiosemicarbazones were prepared from
AB
     H2NCSNHNH2 (I) and the corresponding ketone. PhCH:CHCOPh (10.4 g.), 4.5
     g. I, and 500 cc. EtOH were refluxed 12 hrs., the EtOH distilled off, the
     solution boiled with ligroine, and the residue crystallized from 75% EtOH to
give
     benzalacetophenone thiosemicarbazone (II), m. 143°. The following
     derivative of I and II were prepared (substituent, m.p. of ketone, and m.p. of
     II given): 4-MeO, -, 142° (m. 190° from AcOH); 4-EtO, -,
     157°; 4-PrO, 59°, 164°; 4-allyloxy, 66°,
     159°; 4-BuO, 64° 150°; 4-PhO, 94° 171°;
     3-MeO, -, 146°; 2-MeO, -, 174°; 3,4-di-MeO, -, 165°;
     4,4'-di-MeO, -, 180°; 4-MeO-4'-Me, -, 199°; 4-MeO-4'-Cl, -,
     217°; 4-EtO-4'-Me, 99°, 55° 4-AcO, -, 186°;
     4-HO, -, 244°; 4-AcNH, -, 205°; 4-O2N, -, 185°;
     3,5-di-Br, 133° 227°; 4-MeO-4'-Br, -, 223°; 4-EtSO2,
     145°, 211°; 4-MeO-4'-MeS, 132°, 155°; 4-HO2C, 227°, 253°; 4-HO2C-4'-Cl, 267, 263°; 4'-HO2C,
     224°; 4-MeO-4'-HO2C, -, 224°; 4-β-acrylic acid,
     250°, 258°; 4-HO2CCH2O, 192°, 208°;
     3-HO2CCH2O, 137°, 181°. The following other carbonyl
     compds. were prepared (compound, m.p., m.p. of thiosemicarbazone given):
     \beta-naphthyl (4-methoxystyryl) ketone, -, 117°, \alpha-thienyl
     (4-methoxystyryl)ketone, 87°, 173° and 5-cinnamoyl-8-
     hydroxyquinoline, -, 234°. II and its derivs. are highly effective
     against tubercle bacilli.
=> logoff hold
                                                                    TOTAL
                                                   SINCE FILE
COST IN U.S. DOLLARS
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COST IN U.S. DOLLARS

SINCE FILE TOTAL
ENTRY SESSION
15.96 228.97

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

CA SUBSCRIBER PRICE

SINCE FILE TOTAL
ENTRY SESSION
-3.20 -5.60

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 08:34:51 ON 23 JAN 2008